

Compressive Sensing

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Where Are We?

We ended up with following results

Moral of Story

- *Most signals represented with **finite** numbers of entries*
- *These entries are presented by a vector which is*
*Either **sparse** or **compressible***
- *We could recover the signal perfectly from **few** samples*

Where Are We?

And following questions

- *How much we could reduce the number of samples?*
- *How could we recover the signal efficiently from the samples?*
- *How should we perform the sampling?*
- *Can we guarantee the perfect recovery of a signal?*

Now, we try to answer

Our First Try for Sparse Recovery

Sparse Recovery: *Optimal Approach*

We now want to solve the *sparse recovery problem*:

Sparse Recovery

For given \mathbf{A} and \mathbf{y} , we intend to find vector \mathbf{x} , such that

$$\mathbf{Ax} = \mathbf{y}$$

$$\text{subject to } \|\mathbf{x}\|_0 \leq s$$

for some sparsity s

At this point, we *do not* care about the complexity

Our First Try for Sparse Recovery

Section 1: Optimal Sparse Recovery via ℓ_0 -Norm Minimization

Sparse Recovery: *Optimal Approach*

Assume that

\mathbf{x}^* is the *unique solution* of *sparse recovery problem*

What does this mean?

Consider the *underdetermined* system of equations

$$\mathbf{Ax} = \mathbf{y}$$

Let \mathcal{Z} be the set of all solutions. This means

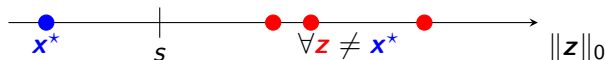
$$\forall \mathbf{z} \in \mathcal{Z} \rightsquigarrow \mathbf{Az} = \mathbf{y} \quad \text{and} \quad \forall \mathbf{z} \notin \mathcal{Z} \rightsquigarrow \mathbf{Az} \neq \mathbf{y}$$

Sparse Recovery: *Optimal Approach*

As a result, we can write

$$\mathbf{x}^* \in \mathcal{Z}$$

What happens if we plot the ℓ_0 -norm of all $\mathbf{z} \in \mathcal{Z}$?



- \mathbf{x}^* is a solution of *sparse recovery problem*

$$\|\mathbf{x}^*\|_0 \leq s$$

- \mathbf{x}^* is the *unique solution*

$$\forall \mathcal{Z} \ni \mathbf{z} \neq \mathbf{x}^* : \|\mathbf{z}\|_0 > s$$

Sparse Recovery: *Optimal Approach*

We can alternatively find the *unique solution* \mathbf{x}^* as

$$\mathbf{x}^* = \operatorname{argmin}_{\mathbf{z} \in \mathcal{Z}} \|\mathbf{z}\|_0$$

From the definition of \mathcal{Z} , we rewrite this optimization as

$$\mathbf{x}^* = \operatorname{argmin} \|\mathbf{z}\|_0 \text{ subject to } \mathbf{A}\mathbf{z} = \mathbf{y}$$

Clearly, \mathbf{x}^* is also the *unique solution* of ℓ_0 -norm minimization

$\mathbf{x}^* \text{ solves sparse recovery} \iff \mathbf{x}^* \text{ solves } \ell_0\text{-norm minimization}$

Sparse Recovery: *Optimal Approach*

ℓ_0 -Norm Recovery

The unique solution of the sparse recovery problem is given by

$$\min \|\mathbf{z}\|_0 \text{ subject to } \mathbf{Az} = \mathbf{y}$$

What does this mean?

*If we have a **unique** solution for sparse recovery problem; then,*

We can recover it for sure via the ℓ_0 -norm recovery

*We hence call ℓ_0 -norm recovery the **optimal** recovery algorithm*

Sparse Recovery: *Optimal Approach*

Now, we know the *optimal* recovery algorithm, but ...

How we could make sure that there *exists* a *unique* solution?

In other words,

What if there exists $\mathbf{x}^\# \neq \mathbf{x}^*$, such that $\mathbf{A}\mathbf{x}^\# = \mathbf{y}$ and

$$\|\mathbf{x}^\#\|_0 \leq s?$$

If $\#$ of samples is larger than some M^* , we can make sure!

We now find out what is M^*

Summary

Up to now, we learned that ...

- The *optimal* approach for sparse recovery is
 ℓ_0 -norm minimization
- We need a minimum number of samples, to make sure
*The sparse recovery is **uniquely** performed*

But what is this minimal number of samples?

*We now find it out for the *optimal* algorithm*

Minimal Number of Samples

Minimal Samples for Sparse Recovery

Depending on the information we have on the sparse signal

$$s \leq M^* \leq 2s$$

We now find out how and when these criteria hold?

We need to first go through some basic definitions

Some Basic Definitions: \mathcal{S} -Subvector

\mathcal{S} -Subvector

Let $\mathbf{x} \in \mathbb{R}^N$, and $\mathcal{S} \subset \{1, \dots, N\}$ of size L , i.e.,

$$\mathcal{S} = \{i_1, \dots, i_L\}$$

where $i_\ell \in \{1, \dots, N\}$ for $\ell \in \{1, \dots, L\}$. The \mathcal{S} -subvector of \mathbf{x} is

$$\mathbf{x}_{\mathcal{S}} = \begin{bmatrix} x_{i_1} \\ \vdots \\ x_{i_L} \end{bmatrix} \in \mathbb{R}^L$$

Some Basic Definitions: \mathcal{S} -Subvector

Example: Consider vector \mathbf{x}

$$\mathbf{x} = \begin{bmatrix} 1 \\ 0 \\ 0 \\ 3 \end{bmatrix}$$

and let $\mathcal{S} = \{2, 4\}$. The \mathcal{S} -subvector of \mathbf{x} is

$$\mathbf{x}_{\mathcal{S}} = \begin{bmatrix} 0 \\ 3 \end{bmatrix}$$

Some Basic Definitions: Column \mathcal{S} -Submatrix

Column \mathcal{S} -Submatrix

Let $\mathbf{A} = [\mathbf{a}_1, \dots, \mathbf{a}_N]$ with $\mathbf{a}_n \in \mathbb{R}^M$, and

$$\mathcal{S} = \{i_1, \dots, i_L\} \subset \{1, \dots, N\}$$

The column \mathcal{S} -submatrix of \mathbf{A} is

$$\mathbf{A}_{\mathcal{S}} = [\mathbf{a}_{i_1}, \dots, \mathbf{a}_{i_L}] \in \mathbb{R}^{M \times L}$$

You can similarly define *row* \mathcal{S} -submatrix

Some Basic Definitions: Column \mathcal{S} -Submatrix

Example: Consider matrix \mathbf{A}

$$\mathbf{A} = \begin{bmatrix} 1 & 0 & 5 & 0 \\ 1 & 0 & 5 & 0 \\ 1 & 0 & 5 & 0 \end{bmatrix}$$

and let $\mathcal{S} = \{2, 3\}$. The column \mathcal{S} -submatrix of \mathbf{A} is

$$\mathbf{A}_{\mathcal{S}} = \begin{bmatrix} 0 & 5 \\ 0 & 5 \\ 0 & 5 \end{bmatrix}$$

Some Basic Definitions: Column \mathcal{S} -Submatrix

Let's do an exercise together which also benefits us later on

Assume that $\text{Supp}(\mathbf{x}) = \mathcal{S}$; show that

$$\mathbf{A}\mathbf{x} = \mathbf{A}_{\mathcal{S}}\mathbf{x}_{\mathcal{S}}$$

We can try it first for

$$\mathbf{x} = \begin{bmatrix} 0 \\ 1 \\ 1 \\ 0 \end{bmatrix} \quad \text{and} \quad \mathbf{A} = [\mathbf{a}_1 \ \mathbf{a}_2 \ \mathbf{a}_3 \ \mathbf{a}_4]$$

Some Basic Definitions: Matrix Kernel

Matrix Kernel

Consider $\mathbf{A} \in \mathbb{R}^{M \times N}$. The *kernel* or *null space* of \mathbf{A} is

$$\ker \mathbf{A} = \left\{ \text{all } \mathbf{x} \in \mathbb{R}^N : \mathbf{A}\mathbf{x} = \mathbf{0} \right\}$$

Two well-known facts about the kernel:

- For *any* matrix \mathbf{A} , the vector of *all zeros* $\mathbf{0} \in \ker \mathbf{A}$
- The kernel of a *fat* matrix *always* has also other elements, i.e.,

$$\ker \mathbf{A} \neq \{\mathbf{0}\}$$

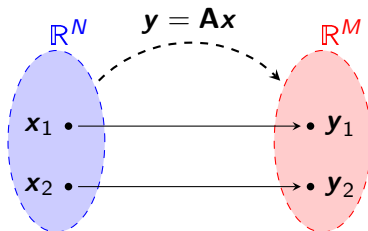
if $\# \text{ rows of } \mathbf{A} < \# \text{ columns } \mathbf{A}$

Some Basic Definitions: Injection

Injective Linear Mapping

Matrix $\mathbf{A} \in \mathbb{R}^{M \times N}$ describes an *injective* linear mapping, if

$$\forall \mathbf{x}_1 \neq \mathbf{x}_2 \in \mathbb{R}^N \rightsquigarrow \mathbf{A}\mathbf{x}_1 \neq \mathbf{A}\mathbf{x}_2$$

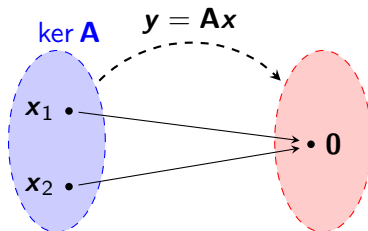


Some Basic Definitions: Injection

Necessary Condition for Injection

Let $\mathbf{A} \in \mathbb{R}^{M \times N}$ describe an *injective* linear mapping; then,

$$\ker \mathbf{A} = \{\mathbf{0}\}$$



Some Basic Definitions: Injection

Main Application of Injection

Let $\mathbf{A} \in \mathbb{R}^{M \times N}$ describe an *injective* linear mapping; then, there exists a matrix $\mathbf{B} \in \mathbb{R}^{N \times M}$

$$\mathbf{BA} = \mathbf{I}_N$$

Roughly speaking,

If \mathbf{A} is *injective*; then, it is *invertible*

Some Basic Definitions: Injection

For an square matrix, *injection* is exactly the *invertibility*

$\mathbf{A} \in \mathbb{R}^{N \times N}$ is *injective*, if it is *invertible*, i.e., $\det \mathbf{A} \neq 0$

In this case, we simply have

$$\mathbf{B} = \mathbf{A}^{-1}$$

Minimal Number of Samples

Section 1: Deriving the Lower Bound

Minimal Number of Samples: *Observation One*

An Initial Lower Bound

Let \mathbf{A} have M rows and \mathbf{x}^* be the *unique* solution of the ℓ_0 -norm minimization. Then, $\|\mathbf{x}^*\|_0 \leq M$

Proof:

Remember the ℓ_0 -norm minimization algorithm

$$\min \|\mathbf{z}\|_0 \text{ subject to } \mathbf{Az} = \mathbf{y}$$

Assume \mathbf{x}^ has $s > M$ nonzero entries. Since \mathbf{x}^* is solution,*

$$\mathbf{Ax}^* = \mathbf{y}$$

Minimal Number of Samples: *Observation One*

An Initial Lower Bound

Let \mathbf{A} have M rows and \mathbf{x}^* be the *unique* solution of the ℓ_0 -norm minimization. Then, $\|\mathbf{x}^*\|_0 \leq M$

Proof:

Let $\mathcal{S} = \text{Supp}(\mathbf{x}^*)$. Then, we can write

$$\mathbf{A}_{\mathcal{S}} \mathbf{x}_{\mathcal{S}}^* = \mathbf{A} \mathbf{x}^* = \mathbf{y}$$

where $\mathbf{A}_{\mathcal{S}} \in \mathbb{R}^{M \times s}$. Since $\mathbf{A}_{\mathcal{S}}$ is fat, we know that

$$\ker \mathbf{A}_{\mathcal{S}} \neq \{\mathbf{0}\}$$

Minimal Number of Samples: *Observation One*

An Initial Lower Bound

Let \mathbf{A} have M rows and \mathbf{x}^* be the *unique* solution of the ℓ_0 -norm minimization. Then, $\|\mathbf{x}^*\|_0 \leq M$

Proof:

Let $\ker \mathbf{A}_S \ni \mathbf{u} \neq \mathbf{0}$, and define $\mathbf{x}^\#$ to be

$\mathbf{x}_S^* + \mathbf{u}$ at entries of S and zero elsewhere $\equiv \mathbf{x}_S^\# = \mathbf{x}_S^* + \mathbf{u}$

Clearly, we have

$$\text{Supp}(\mathbf{x}^\#) \subseteq \text{Supp}(\mathbf{x}^*) \rightsquigarrow \|\mathbf{x}^\#\|_0 \leq \|\mathbf{x}^*\|_0$$

Minimal Number of Samples: *Observation One*

An Initial Lower Bound

Let \mathbf{A} have M rows and \mathbf{x}^* be the *unique* solution of the ℓ_0 -norm minimization. Then, $\|\mathbf{x}^*\|_0 \leq M$

Proof:

Moreover, we can write

$$\begin{aligned}\mathbf{A}\mathbf{x}^\# &= \mathbf{A}_S\mathbf{x}_S^\# = \mathbf{A}_S\mathbf{x}_S^* + \underbrace{\mathbf{A}_S\mathbf{u}}_0 \\ &= \mathbf{A}_S\mathbf{x}_S^* \\ &= \mathbf{y}\end{aligned}$$

Minimal Number of Samples: *Observation One*

An Initial Lower Bound

Let \mathbf{A} have M rows and \mathbf{x}^* be the *unique* solution of the ℓ_0 -norm minimization. Then, $\|\mathbf{x}^*\|_0 \leq M$

Proof:

This means $\mathbf{x}^\#$ is either sparser than, or as sparse as \mathbf{x}^ , i.e.,*

\mathbf{x}^ with $\|\mathbf{x}^*\|_0 > M$ cannot be *unique* solution*

Or equivalently

If \mathbf{x}^ is the unique solution; then, $\|\mathbf{x}^*\|_0 \leq M$*

Minimal Number of Samples: *Observation One*

From observation one, we could conclude that

*The minimal number of samples is **larger than or equal to** s*

In other words,

*With $M < s$ samples, we **cannot** recover a sparse signal*

Minimal Number of Samples

Section 2: Deriving the Upper Bound

Minimal Number of Samples: *Observation Two*

An Upper Bound for Single Sparse Recovery

Let signal $\mathbf{x} \in \mathbb{R}^N$ be s -sparse. Then, there exists a matrix \mathbf{A} with

$$M \geq s + 1$$

rows, such that \mathbf{x} is recovered *uniquely* from samples

$$\mathbf{y} = \mathbf{A}\mathbf{x}$$

via the ℓ_0 -norm minimization algorithm

Proof: See page 53 of the text book

Minimal Number of Samples: *Observation Two*

What does the second finding say?

Let's focus on a *particular* s -sparse signal $\mathbf{x} \in \mathbb{R}^N$; then,

- We design a sampling matrix $\mathbf{A} \in \mathbb{R}^{(s+1) \times N}$ *for signal \mathbf{x}*
- We sample \mathbf{x} as $\mathbf{y} = \mathbf{A}\mathbf{x}$
- We give the samples to ℓ_0 -norm minimization algorithm

We are sure the algorithm recovers *the considered \mathbf{x} uniquely*

Which part of this finding is impractical?

*The sampling matrix **depends on the signal!***

Minimal Number of Samples: *Observation Two*

Could you make it more clear?

Assume we design \mathbf{A} with $s + 1$ rows to sample s -sparse signal \mathbf{x} , such that

\mathbf{x} is recovered uniquely from the samples

Then, there is *no guarantee* that if we use \mathbf{A} to sample another s -sparse signal

$$\mathbf{z} \neq \mathbf{x}$$

we can still uniquely recover \mathbf{z} from the samples $\hat{\mathbf{y}} = \mathbf{A}\mathbf{z}$!

Minimal Number of Samples: *Observation Two*

What kind of sampling matrix are we interested in?

*We are interested in a sampling matrix which
guarantees the **unique** recovery of **all s -sparse signals***

We now try to construct a matrix with this property

Minimal Number of Samples: *Observation Three*

Assume we have *distinct* real numbers $0 < t_1 < \dots < t_N$

$$\mathbf{A} = \begin{bmatrix} t_1^0 & \dots & t_N^0 \\ t_1^1 & \dots & t_N^1 \\ \vdots & & \vdots \\ t_1^{2s-1} & \dots & t_N^{2s-1} \end{bmatrix} = \begin{bmatrix} 1 & \dots & 1 \\ t_1 & \dots & t_N \\ \vdots & & \vdots \\ t_1^{2s-1} & \dots & t_N^{2s-1} \end{bmatrix}$$

For any $S \subset \{1, \dots, N\}$ of size

$$|S| = \# \text{ of elements in } S = 2s$$

the column S -submatrix of \mathbf{A} is *invertible*

See Appendix A of the textbook

Minimal Number of Samples: *Observation Three*

As a result for any S of size $2s$, we can write

$$\mathbf{A}_S \text{ is invertible} \rightsquigarrow \mathbf{A}_S \text{ is injective} \rightsquigarrow \ker \mathbf{A}_S = \{\mathbf{0}\}$$

Now assume that we sample s -sparse signal \mathbf{x} with \mathbf{A}

$$\mathbf{y} = \mathbf{A}\mathbf{x}$$

We want to make sure that there exists *no* \mathbf{x}^\sharp , such that

$$\mathbf{y} = \mathbf{A}\mathbf{x}^\sharp$$

and

$$\|\mathbf{x}^\sharp\|_0 \leq s$$

Deriving the Upper Bound

Minimal Number of Samples: *Observation Three*

Opposite Assumption: Assume there is a \mathbf{x}^\sharp ; then, we have

$$\mathbf{A} \left(\mathbf{x} - \mathbf{x}^\sharp \right) = \mathbf{y} - \mathbf{y} = \mathbf{0}$$

Since $\|\mathbf{x}^\sharp\|_0 \leq \|\mathbf{x}\|_0 \leq s$, we could say

The number of non-zero entries in $\mathbf{x} - \mathbf{x}^\sharp$ is less than $2s$

What is the worst case?

$$\mathbf{x} = \begin{bmatrix} 1 \\ 0 \\ 0 \\ 0 \\ 1 \end{bmatrix} \quad \text{and} \quad \mathbf{x}^\sharp = \begin{bmatrix} 0 \\ 3 \\ 3 \\ 0 \\ 0 \end{bmatrix} \quad \rightsquigarrow \quad \mathbf{x} - \mathbf{x}^\sharp = \begin{bmatrix} 1 \\ -3 \\ -3 \\ 0 \\ 1 \end{bmatrix}$$

Deriving the Upper Bound

Minimal Number of Samples: *Observation Three*

The number of non-zero entries in $\mathbf{x} - \mathbf{x}^\#$ is *at most* $2s$

Therefore, there is a \mathcal{S} of size $2s$ such that

$$\mathbf{A}_{\mathcal{S}} \left(\mathbf{x} - \mathbf{x}^\# \right)_{\mathcal{S}} = \mathbf{A} \left(\mathbf{x} - \mathbf{x}^\# \right) = \mathbf{0}$$

This means that

$$\mathbf{x}_{\mathcal{S}} - \mathbf{x}_{\mathcal{S}}^\# \in \ker \mathbf{A}_{\mathcal{S}}$$

Oops! But, $\ker \mathbf{A}_{\mathcal{S}} = \{\mathbf{0}\}$. This means

$$\mathbf{x} - \mathbf{x}^\# = \mathbf{0} \rightsquigarrow \mathbf{x} = \mathbf{x}^\#$$

Minimal Number of Samples: *Observation Three*

An Upper Bound for Universal Sparse Recovery

There exists a sampling matrix $\mathbf{A} \in \mathbb{R}^{M \times N}$ with

$$M \geq 2s$$

*by which **any** s -sparse signal $\mathbf{x} \in \mathbb{R}^N$ is recovered **uniquely** from*

$$\mathbf{y} = \mathbf{A}\mathbf{x}$$

via the ℓ_0 -norm minimization algorithm

Proof: *We just did it by contradiction! We even know what is \mathbf{A}*

Minimal Number of Samples: *Observation Three*

From observation three, we could conclude that

*The minimal number of samples is **smaller than or equal to** $2s$*

In other words,

*With **$M \geq 2s$** samples, we **definitely** recover an s -sparse signal*

Summary

In a nutshell, we learned that ...

- The *optimal* approach for sparse recovery is
 ℓ_0 -norm minimization
- Minimum samples, we need to do *unique* sparse recovery reads

$$s \leq M^* \leq 2s$$

But can we use these results in practice?

No! We explain why it is so in the next sections

Complexity of ℓ_0 -Norm Minimization

Direct Solution to ℓ_0 -Norm Minimization

ℓ_0 -Norm Minimization

$$\min \|z\|_0 \text{ subject to } \mathbf{A}z = \mathbf{y}$$

What is the direct approach to solve this problem?

We need to find all vectors z that satisfy

$$\mathbf{A}z = \mathbf{y}$$

and then choose the solution with minimum ℓ_0 -norm!

Direct Solution to ℓ_0 -Norm Minimization

ℓ_0 -Norm Minimization

$$\min \|z\|_0 \text{ subject to } \mathbf{A}z = \mathbf{y}$$

What is the direct approach to solve this problem?

*This is however a search with **infinite** feasible points!*

This is not possible to implement!

*This approach seems to be **infeasible***

Direct Solution to ℓ_0 -Norm Minimization

ℓ_0 -Norm Minimization

$$\min \|z\|_0 \text{ subject to } \mathbf{A}z = \mathbf{y}$$

Is there an *alternative* approach?

Well, we could go for the original *sparse recovery problem*, i.e.,

$$\mathbf{A}z = \mathbf{y} \text{ subject to } \|z\|_0 \leq s$$

for some known s

Direct Solution to ℓ_0 -Norm Minimization

ℓ_0 -Norm Minimization: *Alternative Problem*

$$\mathbf{A}\mathbf{z} = \mathbf{y} \text{ subject to } \|\mathbf{z}\|_0 \leq s$$

But how can we solve this *alternative* problem?

We could set $\|\mathbf{z}\|_0 = s$: Consider all possible supports

$$\# \text{ of possible supports} = \binom{N}{s}$$

For each support, we solve $\mathbf{A}\mathbf{z} = \mathbf{y}$

Direct Solution to ℓ_0 -Norm Minimization

ℓ_0 -Norm Minimization: *Alternative Problem*

$$\mathbf{A}\mathbf{z} = \mathbf{y} \text{ subject to } \|\mathbf{z}\|_0 \leq s$$

But how can we solve this *alternative* problem?

Let $\mathcal{S} \subseteq \{1, \dots, N\}$ be a support with $|\mathcal{S}| = s$; then,

$$\mathbf{A}\mathbf{z} = \mathbf{y} \rightsquigarrow \mathbf{A}_{\mathcal{S}}\mathbf{z}_{\mathcal{S}} = \mathbf{y}$$

Depending on $\mathbf{A}_{\mathcal{S}}$, this is *determined*, *under-* or *overdetermined*

Direct Solution to ℓ_0 -Norm Minimization

ℓ_0 -Norm Minimization: *Alternative Problem*

$$\mathbf{A}\mathbf{z} = \mathbf{y} \text{ subject to } \|\mathbf{z}\|_0 \leq s$$

But how can we solve this *alternative* problem?

\mathbf{A}_S has M rows and $s \leq M$ columns

If less than s *linearly independent rows* are in \mathbf{A}_S

Then, the problem is *underdetermined*, but this is *impossible*

Direct Solution to ℓ_0 -Norm Minimization

ℓ_0 -Norm Minimization: *Alternative Problem*

$$\mathbf{A}\mathbf{z} = \mathbf{y} \text{ subject to } \|\mathbf{z}\|_0 \leq s$$

But how can we solve this *alternative* problem?

Why *impossible*? Remember that *underdetermined* means

We have *multiple* solutions to the sparse recovery problem

But we have a *unique* solution. This is hence *impossible*

Direct Solution to ℓ_0 -Norm Minimization

ℓ_0 -Norm Minimization: *Alternative Problem*

$$\mathbf{A}\mathbf{z} = \mathbf{y} \text{ subject to } \|\mathbf{z}\|_0 \leq s$$

But how can we solve this *alternative* problem?

\mathbf{A}_S has M rows and $s \leq M$ columns

With s *independent rows* and no other *conflicting equations*

The problem is *determined* with a *single* solution

Direct Solution to ℓ_0 -Norm Minimization

ℓ_0 -Norm Minimization: *Alternative Problem*

$$\mathbf{A}\mathbf{z} = \mathbf{y} \text{ subject to } \|\mathbf{z}\|_0 \leq s$$

But how can we solve this *alternative* problem?

\mathbf{A}_S has M rows and $s \leq M$ columns

With s *independent rows* and *conflicting equations*

The problem is *overdetermined* with *no* solution

Direct Solution to ℓ_0 -Norm Minimization

ℓ_0 -Norm Minimization: *Alternative Problem*

$$\mathbf{A}\mathbf{z} = \mathbf{y} \text{ subject to } \|\mathbf{z}\|_0 \leq s$$

But how can we solve this *alternative* problem?

Once we are over with $\|\mathbf{z}\|_0 = s$, we go for $\|\mathbf{z}\|_0 = s - 1, \dots$,

Not really needed, since $(s - 1)$ -sparse vector is also s -sparse

We could limit the search to supports with s elements

Direct Solution to ℓ_0 -Norm Minimization

ℓ_0 -Norm Minimization: *Alternative Problem*

$$\mathbf{A}\mathbf{z} = \mathbf{y} \text{ subject to } \|\mathbf{z}\|_0 \leq s$$

But how can we solve this *alternative* problem?

Since the solution is *unique*, we could say

There is only *one* support, for which the problem is *determined*

So, we keep trying till we end up with the *determined* problem

Direct Solution to ℓ_0 -Norm Minimization

How long does it take to find this solution?

Well, it depends!

- If we are *lucky!*
 - We find the solution in *early steps*
- If we are *not lucky!*
 - We find the solution in *final steps*

What does the statistics say?

*We are half of the times *lucky*, half of the time *unlucky**

Direct Solution to ℓ_0 -Norm Minimization

Then, what would be a typical search time?

Well! Let's say we do in average half of the searches:

- We should check C cases, where

$$C = \frac{1}{2} \binom{N}{s}$$

- Say each check takes T sec, we have

$$\text{Search time} = CT = \frac{1}{2} \binom{N}{s} T \text{ sec}$$

Direct Solution to ℓ_0 -Norm Minimization

What would be the value for a typical application?

Consider a typical MRI:

- We have around $N = 10^7$ pixels
- Say the MRI scan is 1% sparse, this mean

$$s = 0.01 \times 10^7 = 10^5$$

- So, we could say

$$C = \frac{1}{2} \binom{N}{s} \geq \frac{1}{2} \left(\frac{N}{s} \right)^s = \frac{1}{2} \times 100^{10^5} = \frac{1}{2} \times 10^{200000}$$

Direct Solution to ℓ_0 -Norm Minimization

What would be the value for a typical application?

Consider a typical MRI:

- Say each search takes only $T = 10^{-10}$ sec; then,

$$\text{Search time} = \frac{1}{2} \times 10^{199990} \text{ sec}$$

- This is more than 10^{199979} *Centenaries!*

Forget about it!

ℓ_0 -norm Minimization is *not a possible* to implement!

NP-hardness of ℓ_0 -Norm Minimization

*One might say: The **direct** approach is not feasible, but*

*How could we conclude that there is **not an efficient way**?!*

Well, as usual

We need to go through some definitions to answer that

Some Definitions: *P-Problem*

Polynomial Computational Complexity

A *polynomial time* algorithm is an algorithm whose number of operations is a polynomial function of input dimension

Example: Finding the minimum entry of a vector

Say the input is a *vector of length N*

- Compare the first two entries and find the minimum
- Compare minimum with next one, and find new minimum
- Keep repeating till the last entry

Some Definitions: *P-Problem*

Polynomial Computational Complexity

A *polynomial time* algorithm is an algorithm whose number of operations is a polynomial function of input dimension

Example: Finding the minimum entry of a vector

We need to do $N - 1$ comparisons in general

$$\# \text{ of operations} = N - 1$$

which is linear in N

Some Definitions: *P-Problem*

P-Problems

*The class of P-problems is the set of all **decision** problems whose solution is given by a polynomial time algorithm*

*Why **decision** problems?*

*Roughly speaking, when it come to **implementation***

*We can reformulate most problem with **decision** problems*

*You can ignore the phrase **decision** for the moment*

Some Definitions: *NP-Problem*

NP-Problems

*The class of nondeterministic polynomial time problems is the set of all decision problems for which there is a polynomial time algorithm to **certify** the solution*

What does that mean?

*Even if we cannot **solve** the problem in polynomial time,*

*When a solution is given, we can **check** it in polynomial time*

Clearly, we can say

$$\{P\text{-Problems}\} \subseteq \{NP\text{-Problems}\}$$

Some Definitions: *NP-Problem*

Example: Exact Cover by Subsets of Three

We are given with a set

$$\{\mathcal{I}_1, \dots, \mathcal{I}_N\}$$

Each \mathcal{I}_n contains three integers from $\{1, \dots, m\}$

We look for J *non-overlapping* subsets $\mathcal{I}_{i_1}, \dots, \mathcal{I}_{i_J}$, such that

$$\mathcal{I}_{i_1} \cup \dots \cup \mathcal{I}_{i_J} = \{1, \dots, m\}$$

Some Definitions: *NP-Problem*

Example: Exact Cover by Subsets of Three

*If someone claims that $\mathcal{I}_{i_1}, \dots, \mathcal{I}_{i_j}$ is a **solution***

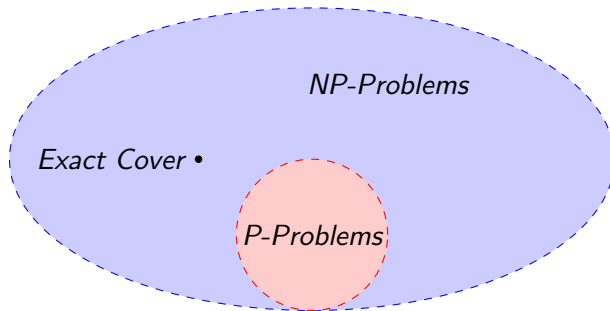
*We can **check** the validity of this claim in polynomial time*

Therefore, we could say

Exact Cover by Subsets of Three $\in \{\text{NP-Problems}\}$

Some Definitions: *NP-Problem*

Where does *Exact Cover by Subsets of Three* locate?



Some Definitions: *NP-Hardness*

NP-Hard Problems

*The class of NP-hard problems is the set of all problems whose solving algorithm is **converted** to the solving algorithm of any NP-problem in polynomial time*

What does that mean?

First of all, note that

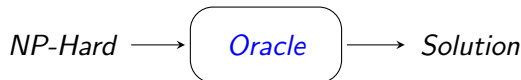
*We **don't** know any polynomial time algorithm
which can solve an NP-hard problem*

Some Definitions: *NP-Hardness*

NP-Hard Problems

*The class of NP-hard problems is the set of all problems whose solving algorithm is **converted** to the solving algorithm of any NP-problem in polynomial time*

*Assume there is an **oracle machine***



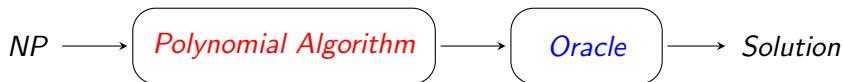
*This machine solves **only one** NP-hard problem*

Some Definitions: *NP-Hardness*

NP-Hard Problems

*The class of NP-hard problems is the set of all problems whose solving algorithm is **converted** to the solving algorithm of any NP-problem in polynomial time*

*Then we can make a **polynomial time** algorithm out of it*



*for **any** NP-problem*

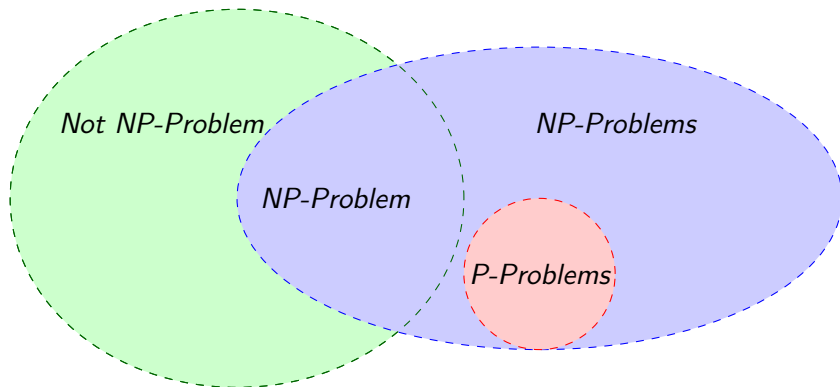
Some Definitions: *NP-Hardness*

We could say

*NP-hard problems are **core** **hard** problems*

*If one day we solve **one**, we solve **all** NP-problems*

Some Definitions: *NP-Hardness*



Some Definitions: *NP-Complete*

*But what if someone **claims** that*

*He/she got a **solution** to an NP-hard problem?*

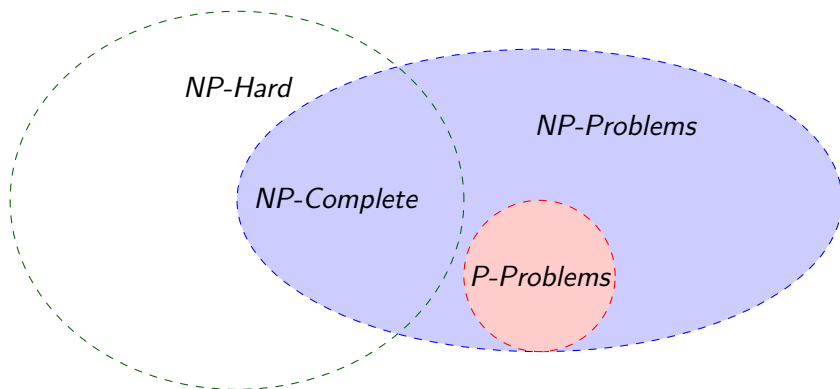
*Of course, we need to be able to **verify** the claim*

*Better if the problem is an **NP-problem***

NP-Complete Problems

*The class of NP-complete problems is the set of all **NP-hard problems** which are **NP-problems** as well*

Some Definitions: *NP-Complete*



Some Definitions: *NP-Complete*

We could say

*NP-complete problems are **atom** **hard** problems*

*If one day we come up with a **solution**,*

- *We can **validate** the solution*
- *We can solve **all NP-problems***

Examples of NP-complete problems

- ***The salesman problem***
- ***Exact cover by subsets of three***

NP-Hardness of Optimal Recovery

What is the use of all of these?

We use these definitions to check whether

*A problem is **feasible** to solve **at the moment** or not*

How can we do it?

*If a given problem describes an **NP-complete** problem*

*Our current tools **cannot** solve the given problem!*

NP-Hardness of Optimal Recovery

Let's get back to *exact cover by subsets of three*

$\mathcal{I}_1, \dots, \mathcal{I}_N$ contain three integers from $\{1, \dots, m\}$

We want to cover $\{1, \dots, m\}$ by J *non-overlapping* of them

Example: Assume $m = 9$ and $N = 4$ with

$$\mathcal{I}_1 = \{1, 2, 3\} , \mathcal{I}_2 = \{3, 4, 5\} , \mathcal{I}_3 = \{4, 5, 6\} , \mathcal{I}_4 = \{7, 8, 9\}$$

Then, the solution has $J = 3$ subsets which are

$$\mathcal{I}_1, \mathcal{I}_3, \mathcal{I}_4$$

NP-Hardness of Optimal Recovery

Let's get back to *exact cover by subsets of three*

$\mathcal{I}_1, \dots, \mathcal{I}_N$ contain three integers from $\{1, \dots, m\}$

We want to cover $\{1, \dots, m\}$ by J *non-overlapping* of them

Let us show each \mathcal{I}_n by a vector $\mathbf{a}_n \in \{0, 1\}^m$, e.g.,

$$\mathcal{I}_1 = \{1, 2, 3\} \rightsquigarrow \mathbf{a}_1 = [1, 1, 1, 0, \dots, 0]^T$$

and construct the matrix \mathbf{A} as

$$\mathbf{A} = [\mathbf{a}_1, \dots, \mathbf{a}_N]$$

NP-Hardness of Optimal Recovery

Let's get back to *exact cover by subsets of three*

$\mathcal{I}_1, \dots, \mathcal{I}_N$ contain three integers from $\{1, \dots, m\}$

We want to cover $\{1, \dots, m\}$ by J *non-overlapping* of them

Now let \mathbf{y} be the vector of m ones

$$\mathbf{y} = \underbrace{[1, \dots, 1]^T}_{m \text{ times}}$$

NP-Hardness of Optimal Recovery

Let's get back to *exact cover by subsets of three*

$\mathcal{I}_1, \dots, \mathcal{I}_N$ contain three integers from $\{1, \dots, m\}$

We want to cover $\{1, \dots, m\}$ by J *non-overlapping* of them

Collection of J subsets could be shown by $\mathbf{x} \in \{0, 1\}^N$, e.g.,

$$\mathcal{I}_1, \mathcal{I}_3, \mathcal{I}_4 \rightsquigarrow \mathbf{x} = \begin{bmatrix} 1 \\ 0 \\ 1 \\ 1 \end{bmatrix}$$

NP-Hardness of Optimal Recovery

Let's get back to *exact cover by subsets of three*

$\mathcal{I}_1, \dots, \mathcal{I}_N$ contain three integers from $\{1, \dots, m\}$

We want to cover $\{1, \dots, m\}$ by J *non-overlapping* of them

Since we aim to *cover* $\{1, \dots, m\}$, we look for an \mathbf{x}

$$\mathbf{Ax} = \mathbf{y}$$

\mathbf{x} is an sparse vector which contains only $J = m/3$ non-zero entry

NP-Hardness of Optimal Recovery

Let's get back to *exact cover by subsets of three*

$\mathcal{I}_1, \dots, \mathcal{I}_N$ contain three integers from $\{1, \dots, m\}$

We want to cover $\{1, \dots, m\}$ by J *non-overlapping* of them

\mathbf{x} is the sparsest possible solution

Any sparser \mathbf{x} would not cover $\{1, \dots, m\}$

We could hence say

$$\mathbf{x} = \operatorname{argmin} \|\mathbf{z}\|_0 \quad \text{subject to } \mathbf{Az} = \mathbf{y}$$

NP-Hardness of Optimal Recovery

What can we now conclude?

ℓ_0 minimization *reduces* to an *NP-complete* problem

Thus, ℓ_0 minimization is an NP-hard problem

*We do **not** know any algorithm which solves it in a *feasible* time*

This is why we go for other algorithms!

Final Points

Summary

- The *optimal* approach for sparse recovery is
 ℓ_0 -norm minimization
- For *unique* sparse recovery via the optimal approach at least

$$s \leq M^* \leq 2s$$

samples should be collected

- ℓ_0 -norm minimization is an *NP-hard* problem
It cannot be implemented in practice

What We Learn Next?

- *We learn sub-optimal algorithms which can be **implemented***
- *Using these algorithms,*

*The **minimum number of samples** changes*

Which Parts of Textbooks?

We are over with this part

I would suggest to go over the text book

*A Mathematical Introduction to Compressive Sensing
S. Foucar and H. Rauhut, Book, 2013*

and study the following part:

- *Chapter 2: Sections 2.2 and 2.3, Pages 48–59*